

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121223\
 Data File : BF136511.D
 Acq On : 12 Dec 2023 12:26
 Operator : CG\JU
 Sample : PB157455BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB157455BS

Quant Time: Dec 12 13:19:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 06 09:40:27 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/14/2023
 Supervised By :mohammad ahmed 12/14/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	128635	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	551530	20.000	ng	0.00
39) Acenaphthene-d10	9.833	164	293123	20.000	ng	0.00
64) Phenanthrene-d10	11.327	188	548616	20.000	ng	0.00
76) Chrysene-d12	13.980	240	281695	20.000	ng	0.00
86) Perylene-d12	15.456	264	262413	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	993542	114.066	ng	0.02
7) Phenol-d6	6.451	99	1241085	114.203	ng	0.00
23) Nitrobenzene-d5	7.363	82	753199	73.828	ng	0.00
42) 2,4,6-Tribromophenol	10.633	330	431064	119.269	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1604289	74.680	ng	0.00
79) Terphenyl-d14	12.921	244	1781837	83.372	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.657	88	116215	33.733	ng	97
3) Pyridine	3.416	79	268362	32.165	ng	96
4) n-Nitrosodimethylamine	3.381	42	138579	38.231	ng	99
6) Aniline	6.463	93	388717	35.394	ng	# 6
8) 2-Chlorophenol	6.586	128	393544	41.387	ng	98
9) Benzaldehyde	6.351	77	199661	35.391	ng	98
10) Phenol	6.463	94	452145	39.100	ng	82
11) bis(2-Chloroethyl)ether	6.539	93	344340	39.949	ng	98
12) 1,3-Dichlorobenzene	6.739	146	397845	37.203	ng	99
13) 1,4-Dichlorobenzene	6.816	146	405471	37.378	ng	98
14) 1,2-Dichlorobenzene	6.969	146	388081	38.044	ng	98
15) Benzyl Alcohol	6.945	79	311446	42.798	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.075	45	424112	38.257	ng	94
17) 2-Methylphenol	7.063	107	317423	41.319	ng	94
18) Hexachloroethane	7.304	117	143014	37.761	ng	98
19) n-Nitroso-di-n-propyla...	7.216	70	252949	40.350	ng	99
20) 3+4-Methylphenols	7.216	107	390363	40.938	ng	# 84
22) Acetophenone	7.210	105	534934	39.277	ng	97
24) Nitrobenzene	7.381	77	378439	36.874	ng	100
25) Isophorone	7.622	82	713416	38.923	ng	98
26) 2-Nitrophenol	7.692	139	210040	41.629	ng	99
27) 2,4-Dimethylphenol	7.739	122	354526	56.985	ng	99
28) bis(2-Chloroethoxy)met...	7.833	93	444218	39.441	ng	98
29) 2,4-Dichlorophenol	7.939	162	335574	40.864	ng	99
30) 1,2,4-Trichlorobenzene	8.016	180	360090	37.605	ng	96
31) Naphthalene	8.098	128	1138941	37.447	ng	99
32) Benzoic acid	7.880	122	268688	43.535	ng	99
33) 4-Chloroaniline	8.151	127	295562	27.258	ng	98
34) Hexachlorobutadiene	8.210	225	207826	36.622	ng	99
35) Caprolactam	8.533	113	116214m	45.955	ng	
36) 4-Chloro-3-methylphenol	8.639	107	366669	42.053	ng	98
37) 2-Methylnaphthalene	8.792	142	752144	38.871	ng	95
38) 1-Methylnaphthalene	8.892	142	704729	37.116	ng	97
40) 1,2,4,5-Tetrachloroben...	8.957	216	368845	40.446	ng	99
41) Hexachlorocyclopentadiene	8.939	237	365768	108.263	ng	99
43) 2,4,6-Trichlorophenol	9.069	196	247916	41.751	ng	97

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44) 2,4,5-Trichlorophenol	9.122	196	268108	40.455	ng	95
46) 1,1'-Biphenyl	9.257	154	984377	39.467	ng	98
47) 2-Chloronaphthalene	9.280	162	737927	38.013	ng	98
48) 2-Nitroaniline	9.380	65	212583	40.830	ng	96
49) Acenaphthylene	9.698	152	1109453	40.119	ng	99
50) Dimethylphthalate	9.563	163	899923	41.700	ng	99
51) 2,6-Dinitrotoluene	9.627	165	199762	41.145	ng	95
52) Acenaphthene	9.869	154	696572	39.308	ng	97
53) 3-Nitroaniline	9.798	138	166880	33.660	ng	95
54) 2,4-Dinitrophenol	9.910	184	218411	88.976	ng	99
55) Dibenzofuran	10.045	168	1018192	39.331	ng	99
56) 4-Nitrophenol	9.974	139	305391	82.169	ng	93
57) 2,4-Dinitrotoluene	10.033	165	261527	40.719	ng	97
58) Fluorene	10.386	166	817806	39.353	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.163	232	231978	43.699	ng	98
60) Diethylphthalate	10.269	149	867968	40.901	ng	99
61) 4-Chlorophenyl-phenyle...	10.380	204	400136	39.498	ng	96
62) 4-Nitroaniline	10.422	138	200983	40.826	ng	99
63) Azobenzene	10.539	77	764439	39.576	ng	96
65) 4,6-Dinitro-2-methylph...	10.445	198	151989	44.734	ng	90
66) n-Nitrosodiphenylamine	10.504	169	741858	42.142	ng	98
67) 4-Bromophenyl-phenylether	10.869	248	262239	41.113	ng	# 91
68) Hexachlorobenzene	10.933	284	298755	40.525	ng	98
69) Atrazine	11.039	200	330078	65.694	ng	97
70) Pentachlorophenol	11.133	266	304016	73.068	ng	99
71) Phenanthrene	11.357	178	1250986	41.486	ng	100
72) Anthracene	11.404	178	1261966	41.550	ng	99
73) Carbazole	11.563	167	1075589	40.012	ng	100
74) Di-n-butylphthalate	11.898	149	1341249	40.700	ng	100
75) Fluoranthene	12.545	202	1195400	37.829	ng	98
77) Benzidine	12.668	184	311466	44.831	ng	99
78) Pyrene	12.774	202	1173090	42.522	ng	98
80) Butylbenzylphthalate	13.398	149	423875	43.161	ng	91
81) Benzo(a)anthracene	13.968	228	810450	41.304	ng	99
82) 3,3'-Dichlorobenzidine	13.933	252	251915	45.790	ng	98
83) Chrysene	14.009	228	771358	39.956	ng	100
84) Bis(2-ethylhexyl)phtha...	13.962	149	517236	42.962	ng	98
85) Di-n-octyl phthalate	14.580	149	770463	43.325	ng	99
87) Indeno(1,2,3-cd)pyrene	16.968	276	854971	46.898	ng	99
88) Benzo(b)fluoranthene	15.027	252	741644	41.336	ng	99
89) Benzo(k)fluoranthene	15.057	252	696519	41.193	ng	100
90) Benzo(a)pyrene	15.398	252	635978	42.922	ng	99
91) Dibenzo(a,h)anthracene	16.986	278	712646	47.716	ng	100
92) Benzo(g,h,i)perylene	17.427	276	707499	46.153	ng	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

